

# 1,2,2,3-Tetrabromopropane

Inchi:	InChI=1S/C3H4Br4/c4-1-3(6,7)2-5/h1-2H2
InchiKey:	SNLFZAHVOKOBOP-UHFFFAOYSA-N
Formula:	C3H4Br4
SMILES:	BrCC(Br)(Br)CBr
Mol. weight [g/mol]:	359.68
CAS:	54268-02-9

## Physical Properties

Property code	Value	Unit	Source
gf	34.50	kJ/mol	Joback Method
hf	-8.68	kJ/mol	Joback Method
hfus	17.25	kJ/mol	Joback Method
hvap	46.72	kJ/mol	Joback Method
log10ws	-3.41		Crippen Method
logp	3.262		Crippen Method
mcvol	123.130	ml/mol	McGowan Method
pc	6818.86	kPa	Joback Method
tb	529.45	K	Joback Method
tc	791.72	K	Joback Method
tf	365.19	K	Joback Method
vc	0.441	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	176.12	J/molxK	529.45	Joback Method
cpg	181.44	J/molxK	573.16	Joback Method
cpg	186.01	J/molxK	616.87	Joback Method
cpg	189.95	J/molxK	660.58	Joback Method
cpg	193.37	J/molxK	704.30	Joback Method
cpg	196.38	J/molxK	748.01	Joback Method
cpg	199.12	J/molxK	791.72	Joback Method
dvisc	0.0021159	Paxs	365.19	Joback Method
dvisc	0.0014518	Paxs	392.57	Joback Method

dvisc	0.0010463	Paxs	419.94	Joback Method
dvisc	0.0007849	Paxs	447.32	Joback Method
dvisc	0.0006086	Paxs	474.70	Joback Method
dvisc	0.0004852	Paxs	502.07	Joback Method
dvisc	0.0003960	Paxs	529.45	Joback Method
hvapt	57.70	kJ/mol	499.00	NIST Webbook

## Sources

<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>
<b>Crippen Method:</b>	<a href="https://www.cheméo.com/doc/models/crippen_log10ws">https://www.cheméo.com/doc/models/crippen_log10ws</a>
<b>Joback Method:</b>	<a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>
<b>McGowan Method:</b>	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>
<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C54268029&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C54268029&amp;Units=SI</a>

## Legend

<b>cpg:</b>	Ideal gas heat capacity
<b>dvisc:</b>	Dynamic viscosity
<b>gf:</b>	Standard Gibbs free energy of formation
<b>hf:</b>	Enthalpy of formation at standard conditions
<b>hfus:</b>	Enthalpy of fusion at standard conditions
<b>hvap:</b>	Enthalpy of vaporization at standard conditions
<b>hvapt:</b>	Enthalpy of vaporization at a given temperature
<b>log10ws:</b>	Log10 of Water solubility in mol/l
<b>logp:</b>	Octanol/Water partition coefficient
<b>mcvol:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature
<b>tf:</b>	Normal melting (fusion) point
<b>vc:</b>	Critical Volume

Latest version available from:

<https://www.cheméo.com/cid/109-259-4/1-2-2-3-Tetrabromopropane.pdf>

Generated by Cheméo on 2024-05-01 23:12:11.869071824 +0000 UTC m=+16894380.789649144.

Cheméo (<https://www.cheméo.com>) is the biggest free database of chemical and physical data for the process industry.